

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040522\
 Data File : BP009696.D
 Acq On : 05 Apr 2022 13:54
 Operator : CG/JU
 Sample : N2249-03DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-2DDL

Quant Time: Apr 05 15:05:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 13:35:48 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	232099	20.000	ng	0.00
21) Naphthalene-d8	10.510	136	929647	20.000	ng	0.00
39) Acenaphthene-d10	14.369	164	597350	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	1278791	20.000	ng	0.00
76) Chrysene-d12	21.257	240	1398552	20.000	ng	0.00
86) Perylene-d12	23.633	264	1467624	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	180652	13.589	ng	0.00
7) Phenol-d6	6.934	99	123633	6.877	ng	0.02
23) Nitrobenzene-d5	8.887	82	307886	17.599	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	218364	22.153	ng	0.00
45) 2-Fluorobiphenyl	12.992	172	765878	17.774	ng	0.00
79) Terphenyl-d14	19.721	244	1210076	15.530	ng	0.00
Target Compounds						
						Qvalue
27) 2,4-Dimethylphenol	9.751	122	65935	5.424	ng	# 93
31) Naphthalene	10.563	128	1379167	28.801	ng	99
37) 2-Methylnaphthalene	12.181	142	227177	6.761	ng	97
38) 1-Methylnaphthalene	12.398	142	218666	6.681	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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